

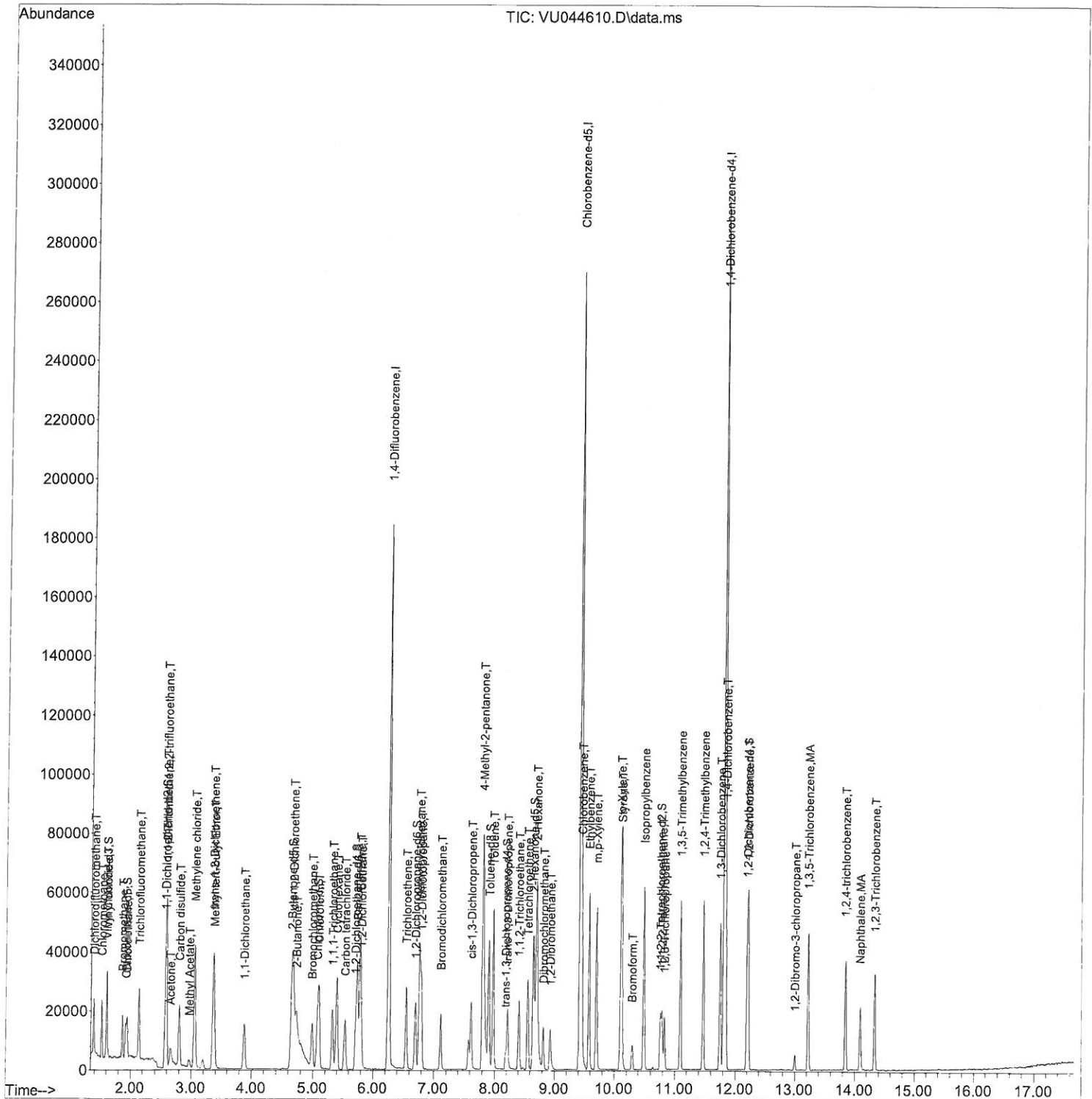
```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090921\  
Data File : VU044610.D  
Acq On    : 09 Sep 2021 13:06  
Operator  : SY/MD  
Sample    : VSTD00122  
Misc      : 25.0mL/MSVOA_U/WATER  
ALS Vial  : 3    Sample Multiplier: 1
```

**Instrument :**  
MSVOA\_U  
**ClientSampleId :**  
VSTD001022

## Manual Integrations

MMDadoda  
9/10/2021 10:50:08 AM

Quant Time: Sep 10 01:06:40 2021  
Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR090921WMA.M  
Quant Title : TRACE VOA SFAM1.0  
QLast Update : Fri Sep 10 01:05:30 2021  
Response via : Initial Calibration



# Quantitation Report (Qedit)

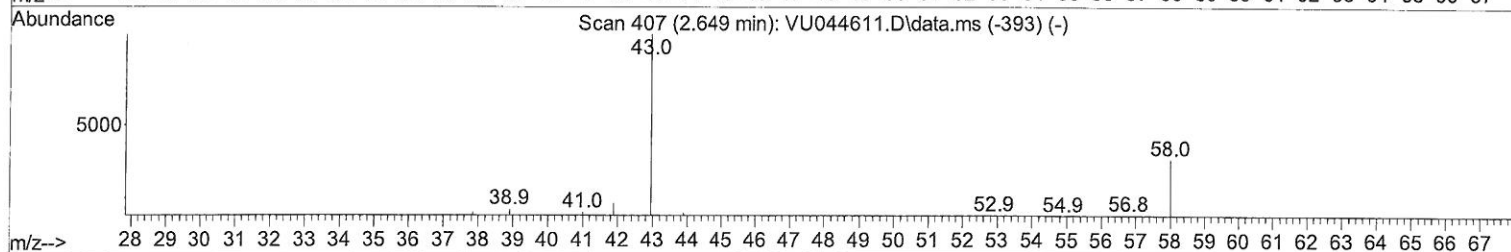
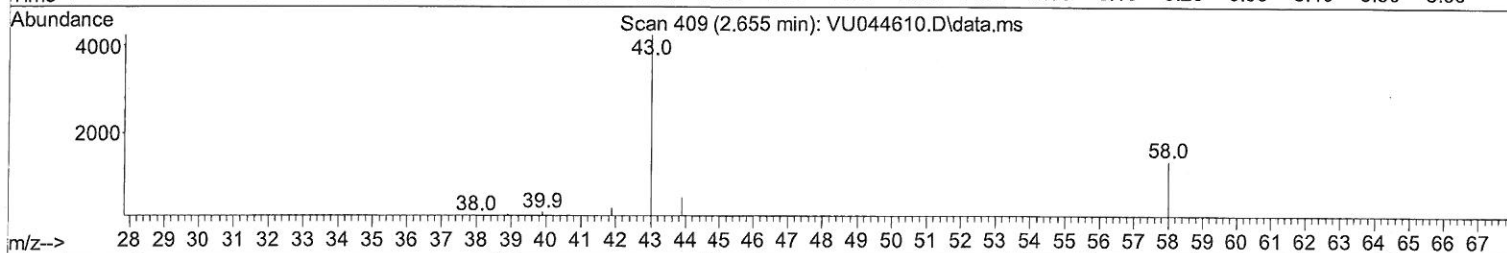
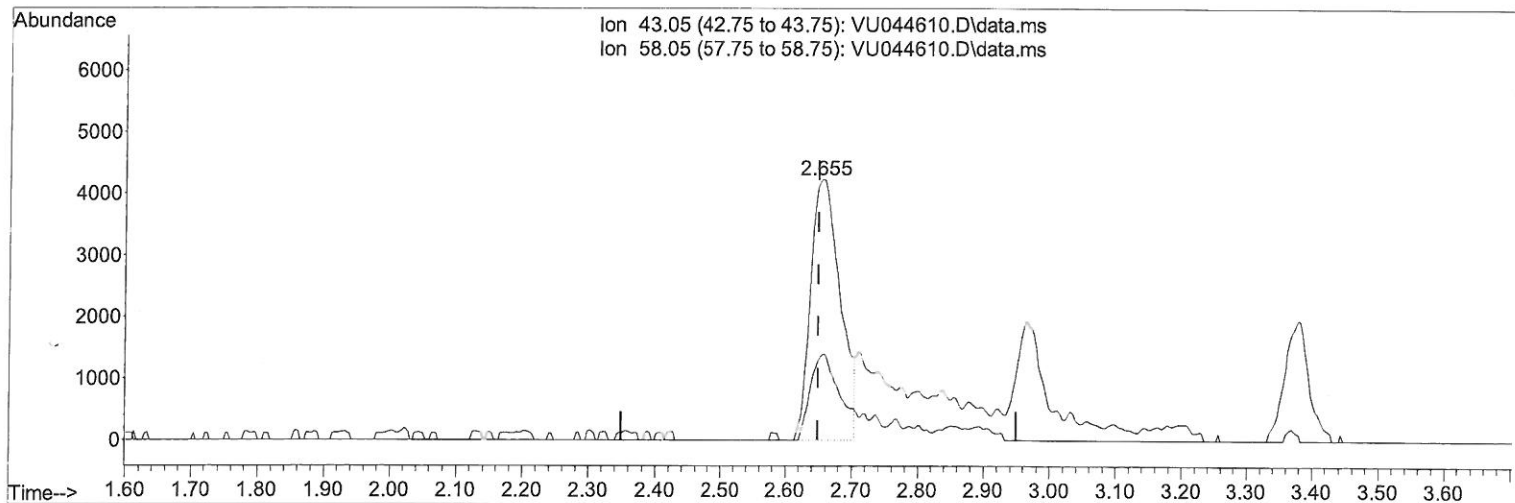
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TIC: VU044610.D\data.ms

(13) Acetone (T)

2.655min (+ 0.006) 5.11 ug/L

response 13033

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 43.05 | 100.00 | 100.00 |
| 58.05 | 21.20  | 33.52  |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

# Quantitation Report (Qedit)

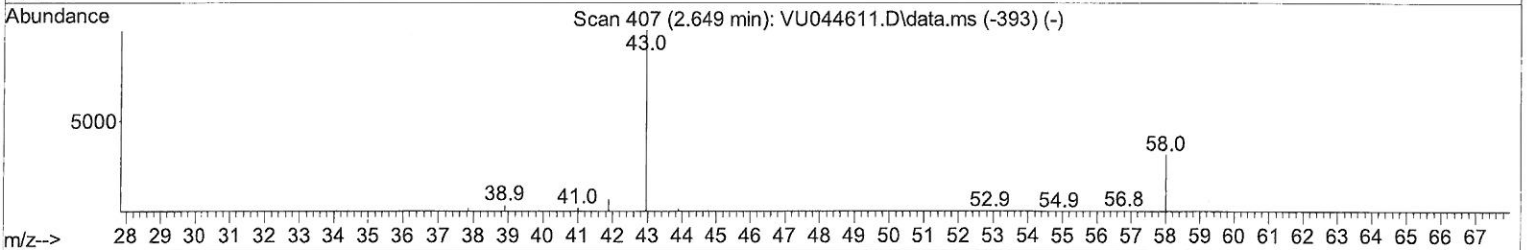
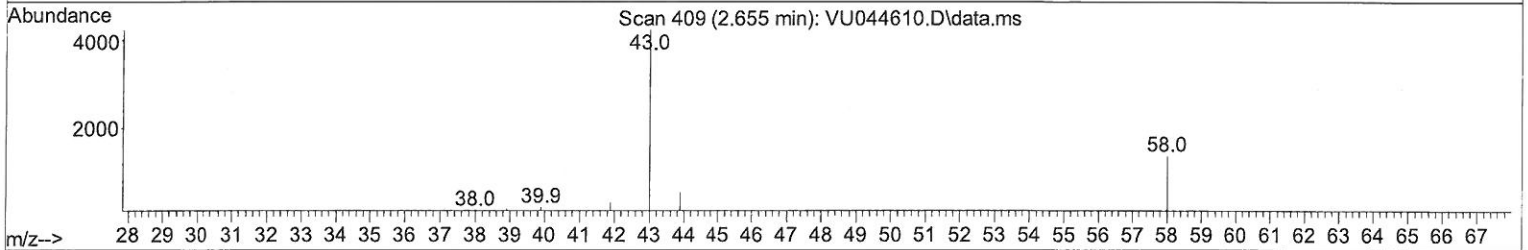
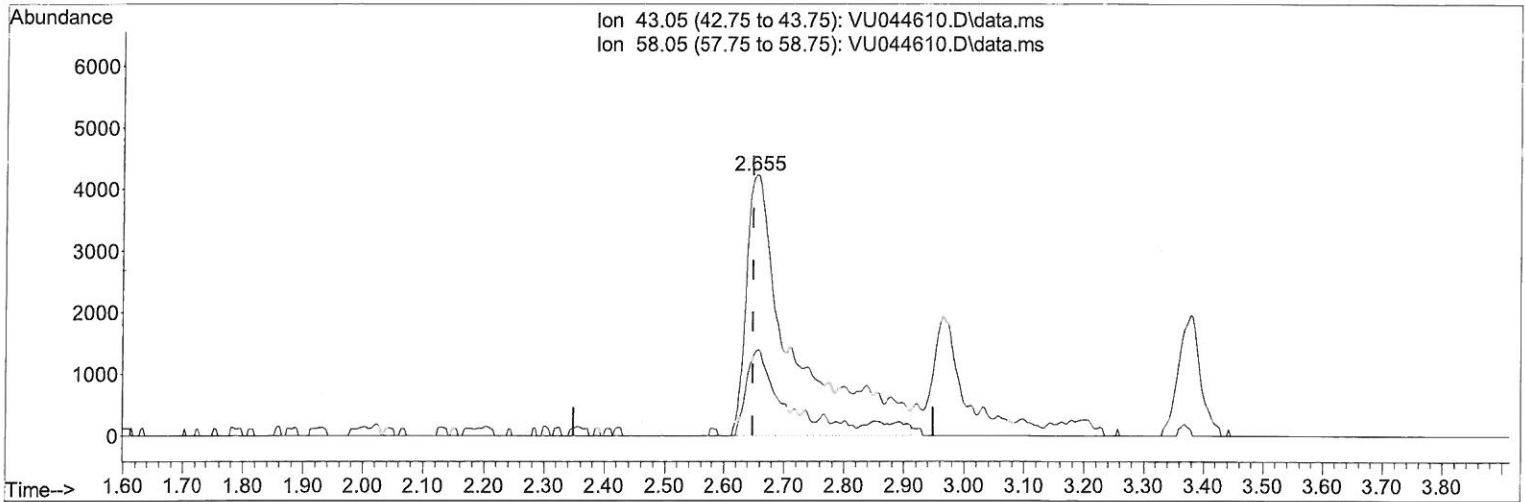
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TIC: VU044610.D\data.ms

(13) Acetone (T)

2.655min (+ 0.006) 8.98 ug/L m

response 22934

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 43.05 | 100.00 | 100.00 |
| 58.05 | 21.20  | 19.05  |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

*MD*  
*9/10/21*

# Quantitation Report (Qedit)

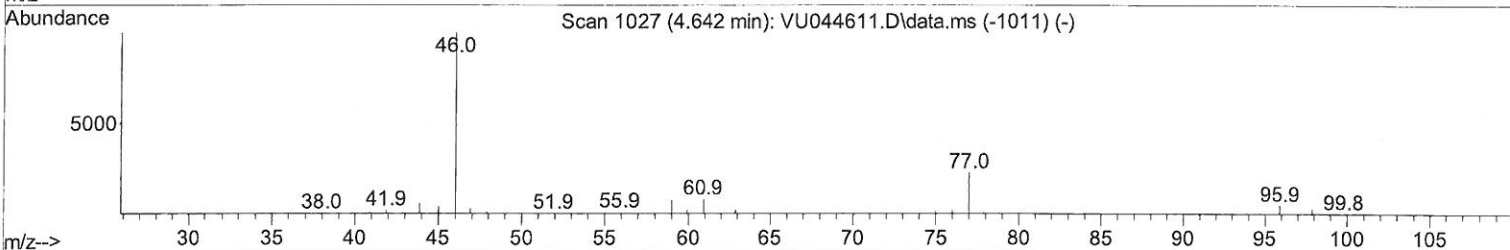
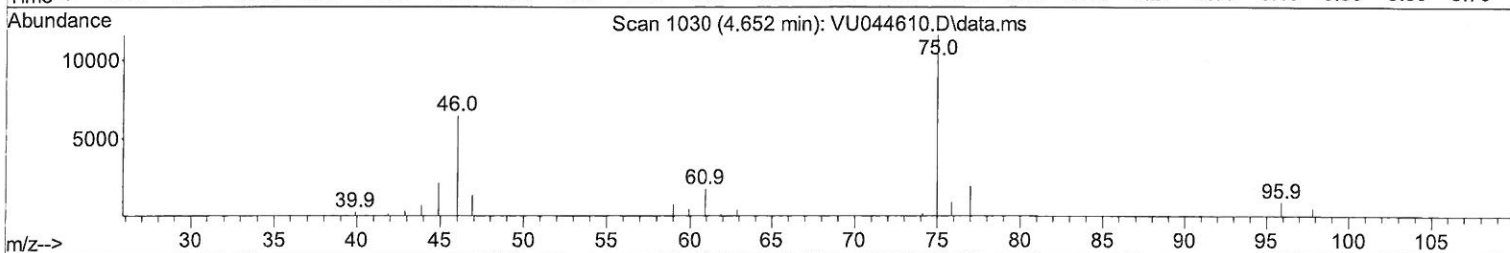
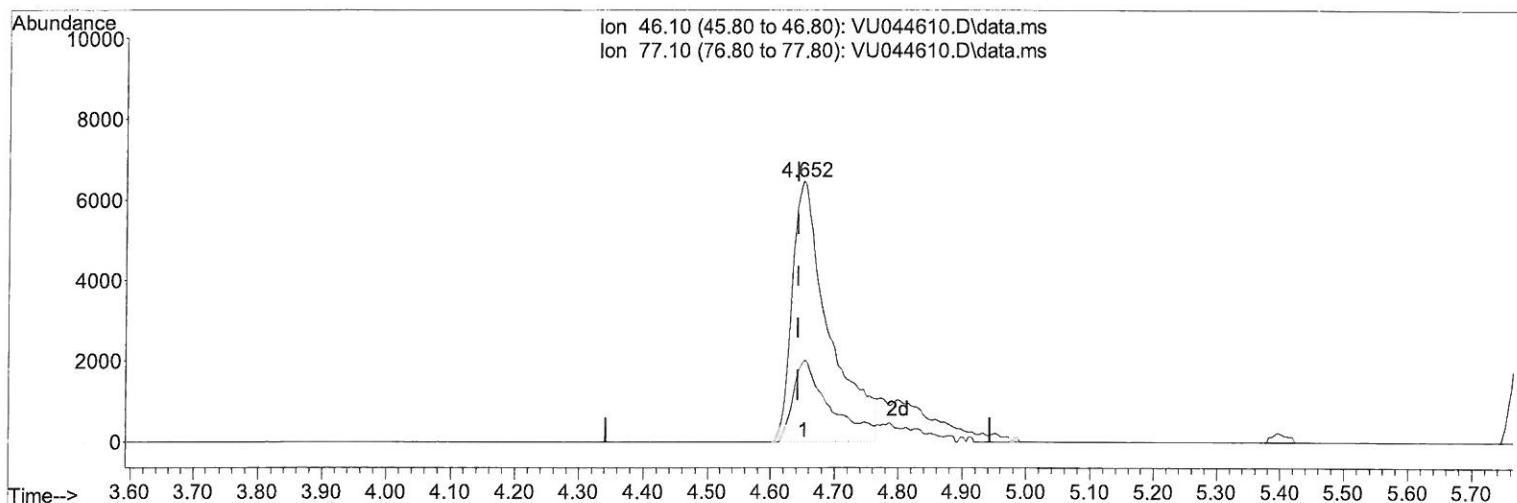
Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU090921\  
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TIC: VU044610.D\data.ms

(20) 2-Butanone-d5 (S)

4.652min (+ 0.009) 8.08 ug/L

response 25801

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 46.10 | 100.00 | 100.00 |
| 77.10 | 17.40  | 29.12# |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |



# Quantitation Report (Qedit)

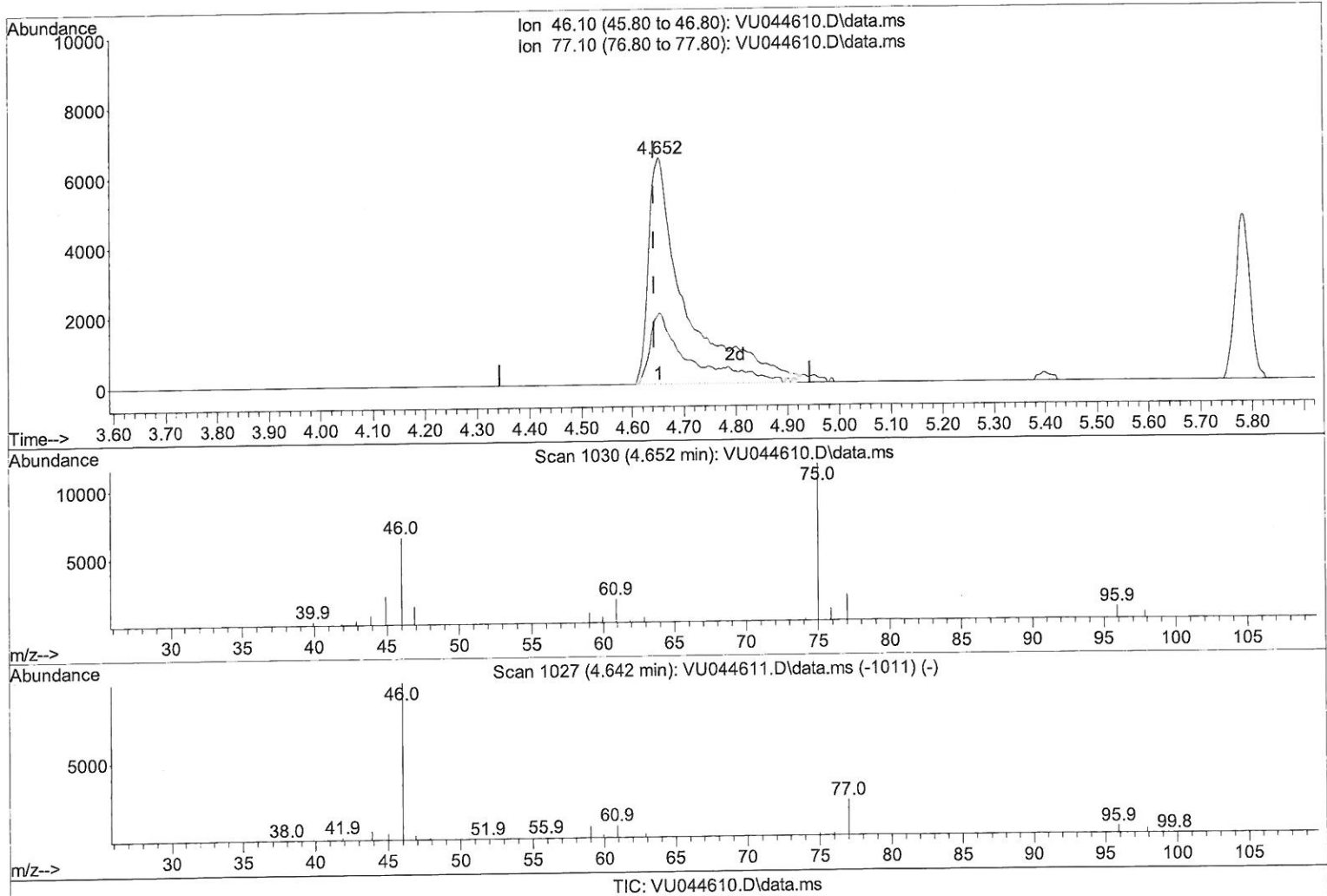
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 Data File : VU044610.D  
 Acq On : 09 Sep 2021 13:06  
 Operator : SY/MD  
 Sample : VSTD00122  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 VSTD001022

Manual Integrations  
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MMDadoda  
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Quant Time: Sep 10 01:06:40 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\SFAMUTR090921WMA.M  
 Quant Title : TRACE VOA SFAM1.0  
 QLast Update : Fri Sep 10 01:05:30 2021  
 Response via : Initial Calibration



(20) 2-Butanone-d5 (S)

4.652min (+ 0.009) 10.02 ug/L m

response 32006

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 46.10 | 100.00 | 100.00 |
| 77.10 | 17.40  | 23.48# |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

MD  
 09/10/21

# Quantitation Report (Qedit)

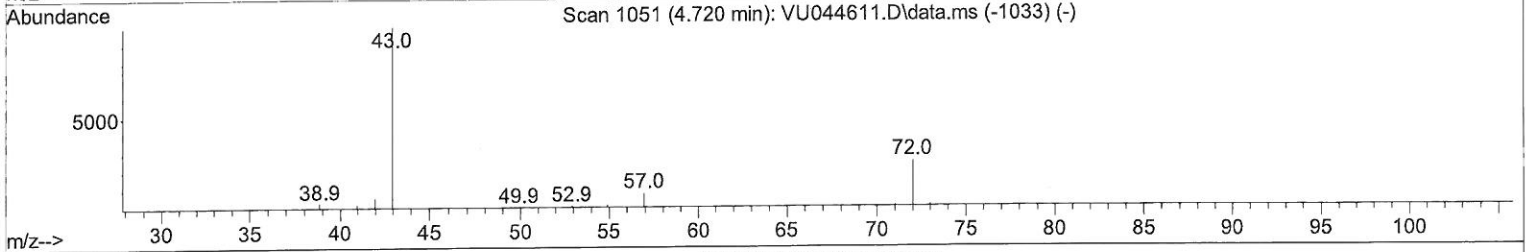
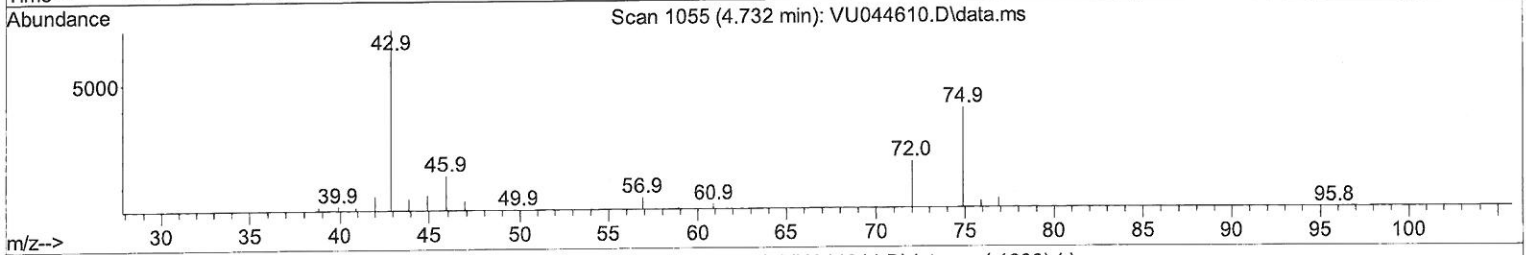
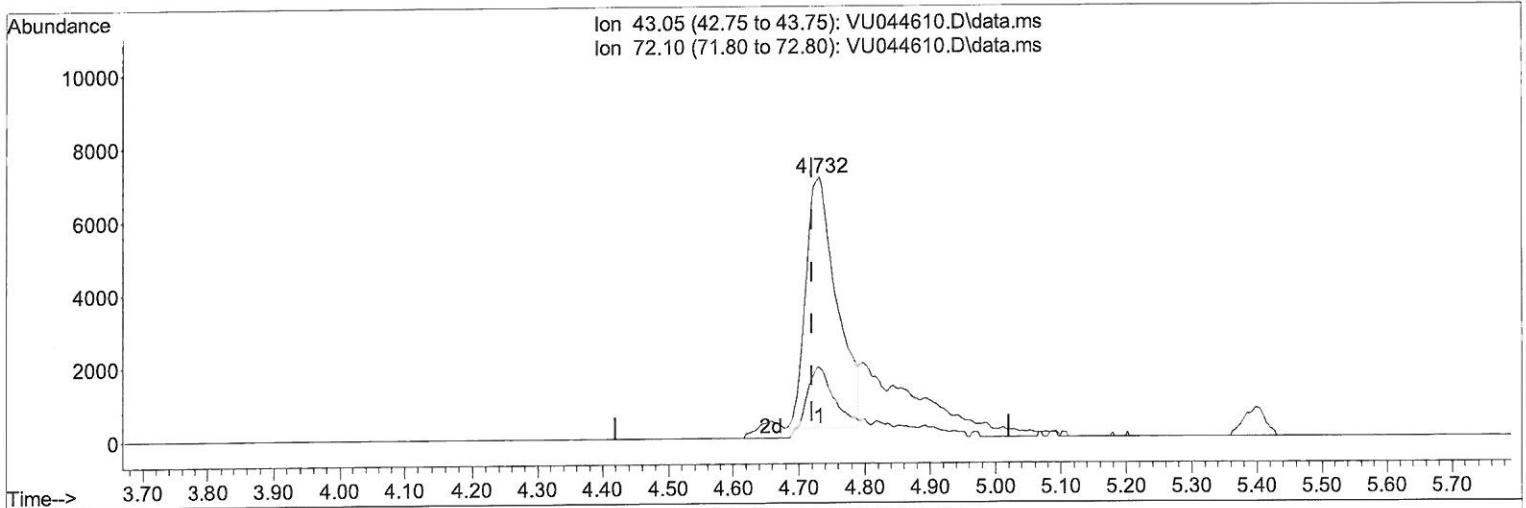
Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU090921\  
 Data File : VU044610.D  
 Acq On : 09 Sep 2021 13:06  
 Operator : SY/MD  
 Sample : VSTD00122  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
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 ClientSampleId :  
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Manual Integrations  
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 Quant Title : TRACE VOA SFAM1.0  
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 Response via : Initial Calibration



TIC: VU044610.D\data.ms

(21) 2-Butanone (T)

4.732min (+ 0.013) 6.03 ug/L

response 22267

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 43.05 | 100.00 | 100.00 |
| 72.10 | 18.30  | 27.83# |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

# Quantitation Report (Qedit)

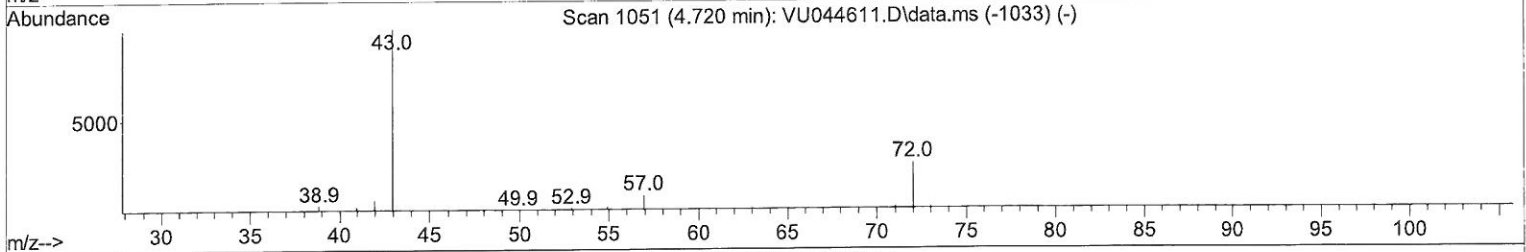
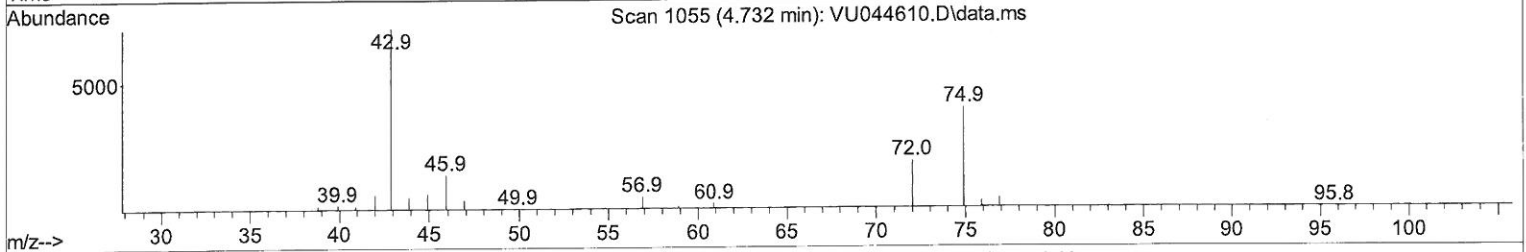
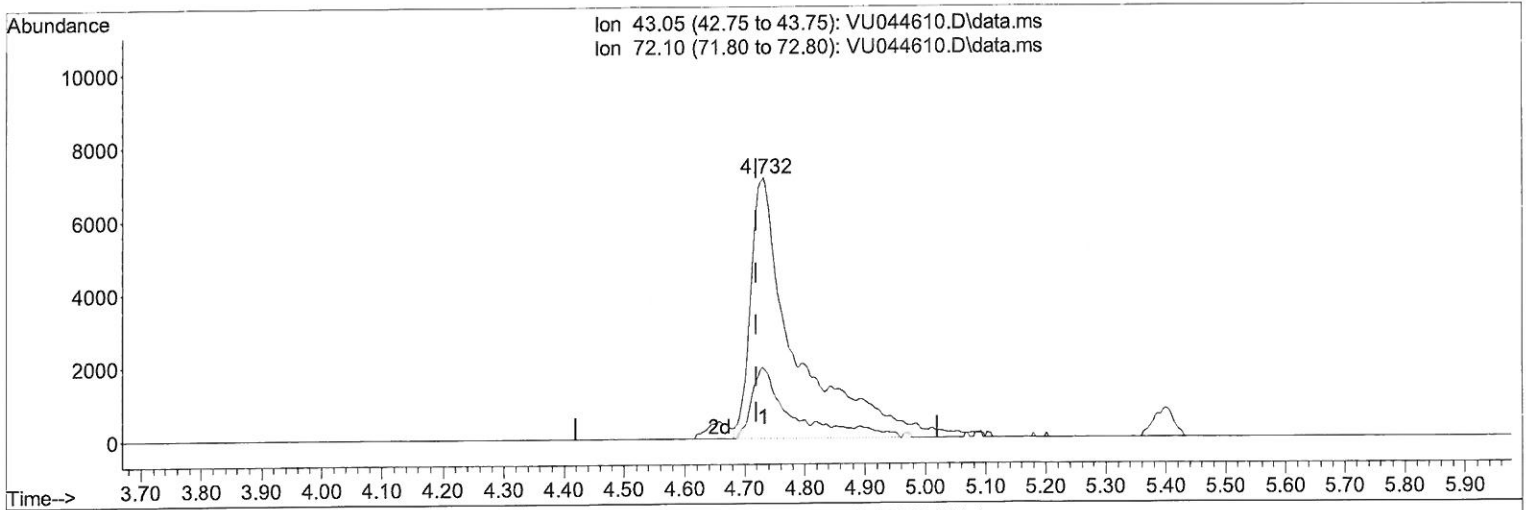
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 Operator : SY/MD  
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 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 VSTD001022

Manual Integrations  
 APPROVED

MMDadoda  
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 Response via : Initial Calibration



TIC: VU044610.D\data.ms

(21) 2-Butanone (T)

4.732min (+ 0.013) 9.71 ug/L m

response 35886

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 43.05 | 100.00 | 100.00 |
| 72.10 | 18.30  | 17.27  |
| 0.00  | 0.00   | 0.00   |
| 0.00  | 0.00   | 0.00   |

7 MD  
 29/10/21

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\U090921\  
 Data File : VU044610.D  
 Acq On : 09 Sep 2021 13:06  
 Operator : SY/MD  
 Sample : VSTD00122  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
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Manual Integrations  
 APPROVED

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 Quant Title : TRACE VOA SFAM1.0  
 QLast Update : Fri Sep 10 01:05:30 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Difluorobenzene        | 6.256  | 114  | 138285   | 5.000  | ug/L   | 0.00     |
| 28) Chlorobenzene-d5          | 9.423  | 117  | 143069   | 5.000  | ug/L   | 0.00     |
| 58) 1,4-Dichlorobenzene-d4    | 11.819 | 152  | 71840    | 5.000  | ug/L   | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 4) Vinyl Chloride-d3          | 1.600  | 65   | 8079     | 0.748  | ug/L   | 0.00     |
| 7) Chloroethane-d5            | 1.919  | 69   | 8201     | 0.857  | ug/L   | 0.00     |
| 11) 1,1-Dichloroethene-d2     | 2.571  | 65   | 3791     | 0.804  | ug/L   | 0.00     |
| 20) 2-Butanone-d5             | 4.652  | 46   | 32006m   | 10.023 | ug/L   | 0.00     |
| 24) Chloroform-d              | 5.073  | 84   | 15990    | 0.826  | ug/L   | 0.00     |
| 26) 1,2-Dichloroethane-d4     | 5.710  | 65   | 9822     | 0.908  | ug/L   | 0.00     |
| 32) Benzene-d6                | 5.735  | 84   | 33017    | 0.819  | ug/L   | 0.00     |
| 36) 1,2-Dichloropropane-d6    | 6.700  | 67   | 10944    | 0.847  | ug/L   | 0.00     |
| 41) Toluene-d8                | 7.906  | 98   | 27345    | 0.766  | ug/L   | 0.00     |
| 43) trans-1,3-Dichloroprop... | 8.185  | 79   | 3433     | 0.741  | ug/L   | 0.00     |
| 46) 2-Hexanone-d5             | 8.648  | 63   | 21780    | 8.432  | ug/L   | 0.00     |
| 56) 1,1,2,2-Tetrachloroeth... | 10.761 | 84   | 9281     | 0.846  | ug/L   | 0.00     |
| 66) 1,2-Dichlorobenzene-d4    | 12.198 | 152  | 10454    | 0.859  | ug/L   | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) Dichlorodifluoromethane    | 1.388  | 85   | 10737    | 0.789  | ug/L   | 99       |
| 3) Chloromethane              | 1.520  | 50   | 13088    | 0.739  | ug/L   | 98       |
| 5) Vinyl chloride             | 1.607  | 62   | 12815    | 0.775  | ug/L   | 96       |
| 6) Bromomethane               | 1.861  | 94   | 6843     | 0.765  | ug/L   | 89       |
| 8) Chloroethane               | 1.938  | 64   | 8248     | 0.806  | ug/L   | 97       |
| 9) Trichlorofluoromethane     | 2.144  | 101  | 14833    | 0.782  | ug/L   | 98       |
| 10) 1,1,2-Trichloro-1,2,2-... | 2.587  | 101  | 8794     | 0.831  | ug/L   | 96       |
| 12) 1,1-Dichloroethene        | 2.587  | 96   | 8413     | 0.812  | ug/L   | 91       |
| 13) Acetone                   | 2.655  | 43   | 22934m   | 8.984  | ug/L   |          |
| 14) Carbon disulfide          | 2.800  | 76   | 26244    | 0.788  | ug/L   | 97       |
| 15) Methyl Acetate            | 2.964  | 43   | 3261     | 0.810  | ug/L   | 94       |
| 16) Methylene chloride        | 3.054  | 84   | 18912    | 1.183  | ug/L   | 96       |
| 17) Methyl tert-butyl Ether   | 3.375  | 73   | 22457    | 0.834  | ug/L   | 100      |
| 18) trans-1,2-Dichloroethene  | 3.366  | 96   | 8777     | 0.812  | ug/L   | 96       |
| 19) 1,1-Dichloroethane        | 3.880  | 63   | 18198    | 0.858  | ug/L   | 98       |
| 21) 2-Butanone                | 4.732  | 43   | 35886m   | 9.713  | ug/L   |          |
| 22) cis-1,2-Dichloroethene    | 4.677  | 96   | 9686     | 0.839  | ug/L   | 95       |
| 23) Bromochloromethane        | 4.983  | 128  | 4352     | 0.867  | ug/L   | 94       |
| 25) Chloroform                | 5.095  | 83   | 19770    | 0.897  | ug/L   | 98       |
| 27) 1,2-Dichloroethane        | 5.803  | 62   | 12782    | 0.887  | ug/L   | 98       |
| 29) 1,1,1-Trichloroethane     | 5.324  | 97   | 14332    | 0.815  | ug/L   | 99       |
| 30) Cyclohexane               | 5.398  | 56   | 15325    | 0.785  | ug/L   | 98       |
| 31) Carbon tetrachloride      | 5.533  | 117  | 11105    | 0.775  | ug/L   | 98       |
| 33) Benzene                   | 5.783  | 78   | 38156    | 0.799  | ug/L   | 100      |
| 34) Trichloroethene           | 6.549  | 95   | 9405     | 0.795  | ug/L   | 96       |
| 35) Methylcyclohexane         | 6.767  | 83   | 14577    | 0.744  | ug/L   | 97       |
| 37) 1,2-Dichloropropane       | 6.796  | 63   | 10596    | 0.844  | ug/L   | 98       |
| 38) Bromodichloromethane      | 7.115  | 83   | 12563    | 0.820  | ug/L   | 95       |
| 39) cis-1,3-Dichloropropene   | 7.616  | 75   | 14061    | 0.798  | ug/L   | 100      |
| 40) 4-Methyl-2-pentanone      | 7.803  | 43   | 76120    | 8.739  | ug/L   | 99       |
| 42) Toluene                   | 7.976  | 91   | 37990    | 0.761  | ug/L   | 99       |



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| Compound                      | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 44) trans-1,3-Dichloropropene | 8.217  | 75   | 11490    | 0.767 | ug/L   | 88       |
| 45) 1,1,2-Trichloroethane     | 8.407  | 97   | 7376     | 0.827 | ug/L   | 95       |
| 47) Tetrachloroethene         | 8.558  | 164  | 6311     | 0.769 | ug/L   | 89       |
| 48) 2-Hexanone                | 8.700  | 43   | 55427    | 8.830 | ug/L   | 96       |
| 49) Dibromochloromethane      | 8.816  | 129  | 7379     | 0.784 | ug/L   | 93       |
| 50) 1,2-Dibromoethane         | 8.931  | 107  | 6832     | 0.801 | ug/L   | 95       |
| 51) Chlorobenzene             | 9.452  | 112  | 24851    | 0.814 | ug/L   | 99       |
| 52) Ethylbenzene              | 9.574  | 91   | 41070    | 0.780 | ug/L   | 100      |
| 53) m,p-Xylene                | 9.700  | 106  | 15637    | 0.789 | ug/L   | 98       |
| 54) o-Xylene                  | 10.105 | 106  | 14790    | 0.771 | ug/L   | 98       |
| 55) Styrene                   | 10.121 | 104  | 24353    | 0.758 | ug/L   | 100      |
| 57) 1,1,2,2-Tetrachloroethane | 10.790 | 83   | 9983     | 0.851 | ug/L   | 98       |
| 59) Bromoform                 | 10.301 | 173  | 3893     | 0.783 | ug/L # | 98       |
| 60) Isopropylbenzene          | 10.491 | 105  | 38656    | 0.754 | ug/L   | 99       |
| 61) 1,2,3-Trichloropropane    | 10.828 | 75   | 7624     | 0.860 | ug/L   | 97       |
| 62) 1,3,5-Trimethylbenzene    | 11.092 | 105  | 30782    | 0.734 | ug/L   | 99       |
| 63) 1,2,4-Trimethylbenzene    | 11.475 | 105  | 31159    | 0.725 | ug/L   | 99       |
| 64) 1,3-Dichlorobenzene       | 11.751 | 146  | 18932    | 0.835 | ug/L   | 98       |
| 65) 1,4-Dichlorobenzene       | 11.841 | 146  | 19894    | 0.865 | ug/L   | 98       |
| 67) 1,2-Dichlorobenzene       | 12.217 | 146  | 17971    | 0.833 | ug/L   | 98       |
| 68) 1,2-Dibromo-3-chloropr... | 13.005 | 75   | 1499     | 0.815 | ug/L   | 88       |
| 69) 1,3,5-Trichlorobenzene    | 13.224 | 180  | 13688    | 0.847 | ug/L   | 96       |
| 70) 1,2,4-trichlorobenzene    | 13.844 | 180  | 10666    | 0.834 | ug/L   | 99       |
| 71) Naphthalene               | 14.092 | 128  | 17381    | 0.718 | ug/L   | 98       |
| 72) 1,2,3-Trichlorobenzene    | 14.336 | 180  | 9357     | 0.800 | ug/L   | 99       |

(#) = qualifier out of range (m) = manual integration (+) = signals summed